



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 06:05 am BST

PDB ID : 5XV1  
Title : Crystal structure of ATG101-ATG13HORMA  
Authors : Kim, B.-W.; Song, H.K.  
Deposited on : 2017-06-26  
Resolution : 2.51 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

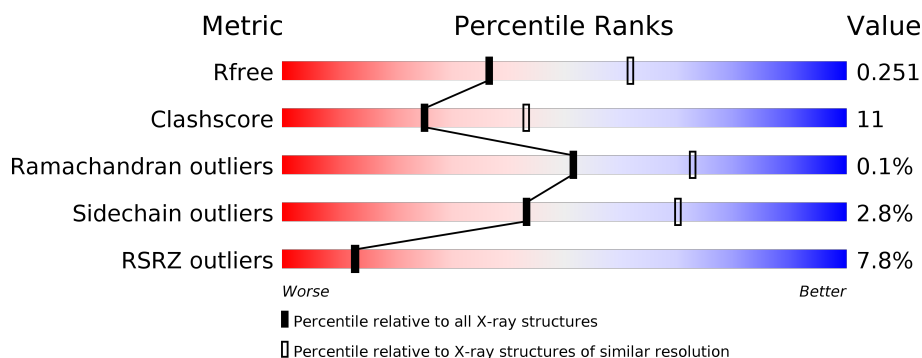
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.51 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	190	<div> <div>7%</div> <div>76%</div> <div>17%</div> <div>5%</div> </div>
1	C	190	<div> <div>7%</div> <div>79%</div> <div>16%</div> <div>•</div> </div>
2	B	218	<div> <div>5%</div> <div>72%</div> <div>24%</div> <div>•</div> </div>
2	D	218	<div> <div>11%</div> <div>73%</div> <div>24%</div> <div>••</div> </div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6095 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Autophagy-related protein 13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	S	0	0	0
			1360	865	232	255	8			
1	C	185	Total	C	N	O	S	0	0	0
			1398	891	239	260	8			

- Molecule 2 is a protein called Autophagy-related protein 101.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	211	Total	C	N	O	S	0	0	0
			1639	1037	279	313	10			
2	D	216	Total	C	N	O	S	0	0	0
			1666	1051	283	322	10			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	40	ALA	LYS	engineered mutation	UNP Q9BSB4
B	41	ALA	LYS	engineered mutation	UNP Q9BSB4
B	42	ALA	GLU	engineered mutation	UNP Q9BSB4
D	40	ALA	LYS	engineered mutation	UNP Q9BSB4
D	41	ALA	LYS	engineered mutation	UNP Q9BSB4
D	42	ALA	GLU	engineered mutation	UNP Q9BSB4

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		
3	B	7	Total	O	0	0
			7	7		
3	C	5	Total	O	0	0
			5	5		

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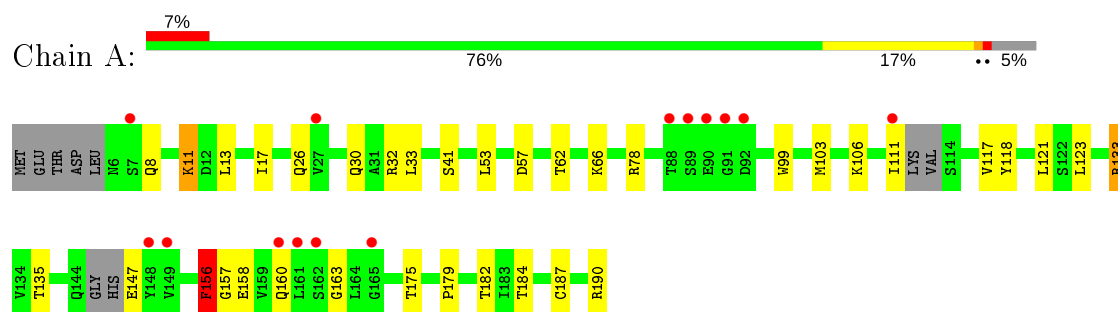
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	11	Total	O	0	0
			11	11		

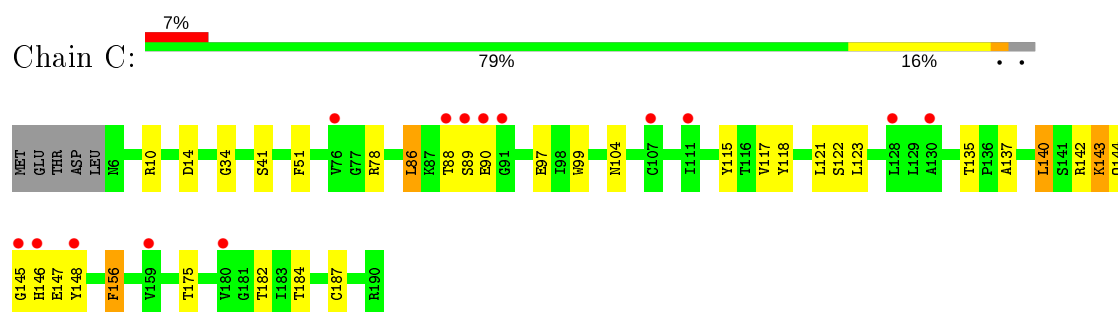
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

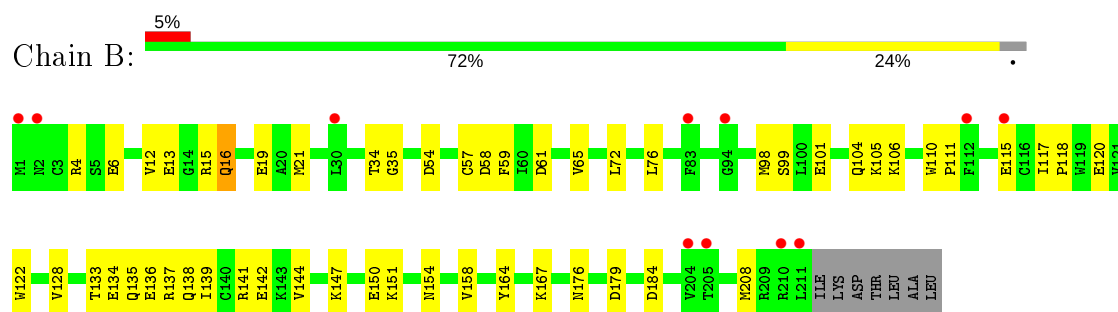
#### • Molecule 1: Autophagy-related protein 13



#### • Molecule 1: Autophagy-related protein 13

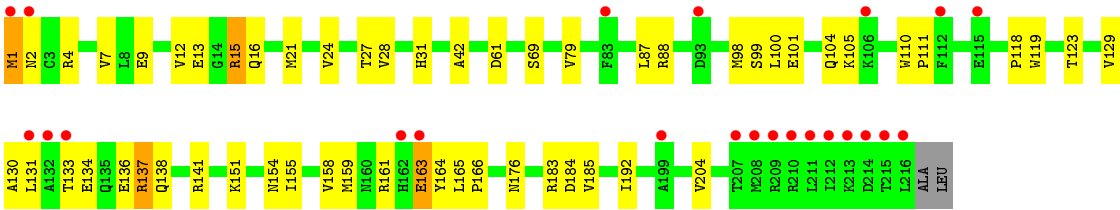


#### • Molecule 2: Autophagy-related protein 101



#### • Molecule 2: Autophagy-related protein 101





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.48Å 125.50Å 101.04Å 90.00° 101.66° 90.00°	Depositor
Resolution (Å)	32.98 – 2.51 38.85 – 2.51	Depositor EDS
% Data completeness (in resolution range)	99.1 (32.98-2.51) 91.3 (38.85-2.51)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.63 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.10.1 _2155	Depositor
R, $R_{free}$	0.210 , 0.251 0.210 , 0.251	Depositor DCC
$R_{free}$ test set	2000 reflections (5.45%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.9	Xtriage
Anisotropy	0.630	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 54.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6095	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.50% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.42	1/1379 (0.1%)	0.60	0/1867
1	C	0.40	0/1421	0.62	1/1926 (0.1%)
2	B	0.39	0/1669	0.56	0/2263
2	D	0.42	0/1696	0.66	3/2300 (0.1%)
All	All	0.41	1/6165 (0.0%)	0.61	4/8356 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	156	PHE	CE1-CZ	-5.10	1.27	1.37

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	137	ARG	NE-CZ-NH2	-8.74	115.93	120.30
2	D	15	ARG	NE-CZ-NH2	-7.67	116.46	120.30
1	C	140	LEU	CA-CB-CG	6.20	129.55	115.30
2	D	137	ARG	NH1-CZ-NH2	6.14	126.16	119.40

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1360	0	1345	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1398	0	1389	23	0
2	B	1639	0	1587	40	0
2	D	1666	0	1591	45	0
3	A	9	0	0	4	0
3	B	7	0	0	1	0
3	C	5	0	0	0	0
3	D	11	0	0	3	0
All	All	6095	0	5912	128	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (128) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:135:GLN:O	2:B:139:ILE:HD12	1.38	1.18
1:A:57:ASP:OD2	3:A:201:HOH:O	1.86	0.93
2:B:15:ARG:HH21	2:B:59:PHE:HD2	1.17	0.87
2:B:58:ASP:O	2:B:141:ARG:NH1	2.12	0.82
2:B:135:GLN:O	2:B:139:ILE:CD1	2.28	0.79
2:B:135:GLN:HG2	2:B:139:ILE:CD1	2.17	0.75
2:D:61:ASP:OD2	2:D:141:ARG:NH1	2.21	0.73
2:D:158:VAL:O	3:D:301:HOH:O	2.06	0.72
1:A:53:LEU:O	3:A:202:HOH:O	2.07	0.71
2:D:4:ARG:CZ	2:D:163:GLU:HG2	2.22	0.70
1:A:135:THR:O	3:A:203:HOH:O	2.10	0.70
1:C:10:ARG:NH1	1:C:14:ASP:OD1	2.26	0.69
2:D:13:GLU:HG3	2:D:131:LEU:HD13	1.75	0.69
1:C:144:GLN:HA	1:C:148:TYR:CE1	2.29	0.68
2:B:13:GLU:OE2	2:B:137:ARG:NH1	2.28	0.67
2:B:179:ASP:OD2	3:B:301:HOH:O	2.13	0.65
2:B:135:GLN:HG2	2:B:139:ILE:HD11	1.77	0.65
2:D:101:GLU:OE1	2:D:118:PRO:HB3	1.97	0.64
1:A:133:ARG:NH1	2:B:54:ASP:OD1	2.30	0.64
2:D:1:MET:N	2:D:2:ASN:HB3	2.12	0.64
1:C:97:GLU:OE1	1:C:135:THR:HG21	1.98	0.64
2:D:161:ARG:HB3	2:D:163:GLU:OE2	1.99	0.63
2:D:4:ARG:CZ	2:D:163:GLU:CG	2.76	0.62
1:C:140:LEU:HD12	1:C:143:LYS:HE3	1.82	0.61
2:B:21:MET:HE1	2:B:98:MET:HB2	1.84	0.60
2:B:15:ARG:NH2	2:B:59:PHE:HD2	1.95	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:133:THR:N	2:D:136:GLU:OE2	2.33	0.59
1:A:175:THR:HG22	1:A:184:THR:HG23	1.83	0.59
2:D:105:LYS:HB2	2:D:176:ASN:HA	1.84	0.59
1:C:86:LEU:HD21	1:C:148:TYR:HB2	1.84	0.59
2:D:79:VAL:HG21	2:D:192:ILE:HD11	1.85	0.58
1:A:8:GLN:HA	1:A:11:LYS:HD2	1.84	0.58
2:B:61:ASP:OD1	2:B:141:ARG:NH2	2.37	0.57
2:B:135:GLN:CG	2:B:139:ILE:HD11	2.36	0.56
2:B:101:GLU:OE1	2:B:118:PRO:HG3	2.05	0.56
1:C:78:ARG:HG2	1:C:156:PHE:CE2	2.41	0.56
1:C:99:TRP:CE3	1:C:187:CYS:HB2	2.41	0.56
1:A:13:LEU:O	1:A:17:ILE:HG13	2.07	0.54
1:A:99:TRP:CE3	1:A:187:CYS:HB2	2.42	0.54
1:A:111:ILE:HG21	1:A:179:PRO:HD3	1.90	0.54
1:C:135:THR:HG22	1:C:137:ALA:H	1.73	0.54
1:A:32:ARG:NH2	1:A:135:THR:HG21	2.22	0.53
2:B:133:THR:HG23	2:B:136:GLU:H	1.73	0.53
2:D:79:VAL:HG11	2:D:192:ILE:HD11	1.88	0.53
2:D:24:VAL:O	2:D:28:VAL:HG23	2.08	0.53
2:D:12:VAL:HG13	2:D:16:GLN:HB2	1.91	0.53
2:B:4:ARG:O	2:B:120:GLU:HA	2.09	0.52
2:B:104:GLN:NE2	2:B:164:TYR:OH	2.41	0.52
1:C:34:GLY:HA2	1:C:142:ARG:HA	1.92	0.52
2:D:13:GLU:OE1	2:D:137:ARG:NE	2.42	0.52
2:B:34:THR:OG1	2:B:35:GLY:N	2.43	0.52
2:B:12:VAL:HB	2:B:16:GLN:HG3	1.91	0.51
1:C:86:LEU:CD2	1:C:148:TYR:HD2	2.24	0.51
2:B:6:GLU:HG3	2:B:158:VAL:HG11	1.93	0.50
2:D:161:ARG:CB	2:D:163:GLU:OE2	2.59	0.50
1:C:51:PHE:HB3	1:C:122:SER:HB2	1.93	0.50
2:D:99:SER:O	2:D:192:ILE:HA	2.11	0.50
2:D:104:GLN:HG3	2:D:119:TRP:CD1	2.46	0.50
2:D:129:VAL:O	3:D:302:HOH:O	2.19	0.50
2:D:161:ARG:C	2:D:163:GLU:OE2	2.51	0.50
2:D:154:ASN:O	2:D:158:VAL:HG13	2.11	0.50
1:A:41:SER:HB2	2:B:184:ASP:HB3	1.94	0.49
1:C:86:LEU:HD23	1:C:148:TYR:HD2	1.77	0.49
2:D:15:ARG:HE	2:D:137:ARG:NH2	2.10	0.49
1:A:26:GLN:O	1:A:30:GLN:HB2	2.13	0.49
1:A:62:THR:HG22	1:A:66:LYS:HD2	1.95	0.49
2:D:104:GLN:CD	2:D:166:PRO:HB3	2.33	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:140:LEU:HD12	1:C:143:LYS:CE	2.43	0.48
2:B:150:GLU:O	2:B:154:ASN:ND2	2.43	0.48
2:B:167:LYS:HD3	2:B:167:LYS:N	2.28	0.47
2:B:72:LEU:O	2:B:76:LEU:HG	2.13	0.47
1:C:117:VAL:O	1:C:121:LEU:HG	2.13	0.47
2:D:100:LEU:HA	2:D:192:ILE:HG22	1.95	0.47
2:B:59:PHE:O	2:B:141:ARG:HG2	2.14	0.47
2:D:163:GLU:OE2	2:D:163:GLU:N	2.48	0.47
2:B:117:ILE:O	2:B:117:ILE:HG13	2.15	0.46
2:B:138:GLN:NE2	2:B:142:GLU:OE2	2.48	0.46
2:B:15:ARG:HE	2:B:15:ARG:HB3	1.54	0.46
2:D:27:THR:O	2:D:31:HIS:HD2	1.99	0.46
1:A:157:GLY:HA3	1:A:158:GLU:C	2.36	0.46
1:A:30:GLN:HA	1:A:33:LEU:CD2	2.46	0.46
2:D:104:GLN:OE1	2:D:164:TYR:OH	2.33	0.46
2:D:4:ARG:CZ	2:D:163:GLU:HG3	2.46	0.46
2:B:134:GLU:O	2:B:137:ARG:HB3	2.16	0.45
2:D:13:GLU:HG2	3:D:302:HOH:O	2.16	0.45
1:C:146:HIS:N	1:C:147:GLU:OE2	2.50	0.45
2:D:7:VAL:HA	2:D:123:THR:O	2.17	0.45
2:D:110:TRP:CG	2:D:111:PRO:HA	2.52	0.45
1:A:157:GLY:HA3	1:A:158:GLU:O	2.16	0.45
1:C:115:TYR:HA	1:C:118:TYR:HB3	1.99	0.45
1:C:175:THR:HG22	1:C:184:THR:HG23	1.98	0.45
2:B:15:ARG:NH1	2:B:137:ARG:CZ	2.80	0.44
1:A:133:ARG:HD3	2:B:65:VAL:CG2	2.48	0.44
1:C:88:THR:HG23	1:C:90:GLU:H	1.82	0.44
1:A:118:TYR:CZ	2:D:42:ALA:HB2	2.53	0.44
1:C:147:GLU:N	1:C:147:GLU:OE2	2.41	0.44
1:C:88:THR:OG1	1:C:89:SER:N	2.50	0.44
1:A:117:VAL:O	1:A:121:LEU:HG	2.18	0.44
1:A:30:GLN:HA	1:A:33:LEU:HD21	2.00	0.44
2:D:155:ILE:O	2:D:159:MET:HB2	2.18	0.43
2:B:12:VAL:O	2:B:128:VAL:HA	2.19	0.43
2:D:9:GLU:O	2:D:151:LYS:HE2	2.18	0.43
2:B:105:LYS:O	2:B:176:ASN:HB3	2.19	0.43
2:B:19:GLU:OE1	2:B:57:CYS:HB3	2.19	0.43
2:B:21:MET:HE2	2:B:98:MET:SD	2.59	0.42
1:C:145:GLY:HA2	1:C:146:HIS:HA	1.60	0.42
1:A:156:PHE:HD1	1:A:156:PHE:HA	1.71	0.42
2:B:16:GLN:OE1	2:B:144:VAL:HG21	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:87:LEU:HD23	2:D:87:LEU:HA	1.92	0.42
1:A:163:GLY:O	3:A:204:HOH:O	2.21	0.42
1:A:147:GLU:OE1	1:A:147:GLU:N	2.52	0.42
2:B:110:TRP:CG	2:B:111:PRO:HA	2.54	0.42
2:B:147:LYS:O	2:B:151:LYS:HG2	2.19	0.41
1:C:104:ASN:ND2	1:C:182:THR:HB	2.35	0.41
2:D:21:MET:HE1	2:D:98:MET:HB2	2.01	0.41
2:D:163:GLU:CD	2:D:163:GLU:H	2.23	0.41
2:D:69:SER:HB2	2:D:185:VAL:HG23	2.02	0.41
2:B:99:SER:HA	2:B:122:TRP:O	2.21	0.41
1:A:78:ARG:HA	1:A:157:GLY:O	2.21	0.41
2:D:130:ALA:C	2:D:131:LEU:HD12	2.40	0.41
1:C:41:SER:HB2	2:D:184:ASP:HB3	2.03	0.40
1:A:106:LYS:O	1:A:182:THR:HG23	2.22	0.40
2:B:106:LYS:HB3	2:B:115:GLU:HG3	2.03	0.40
2:D:4:ARG:NE	2:D:163:GLU:HG2	2.35	0.40
2:D:183:ARG:HG2	2:D:184:ASP:OD1	2.21	0.40
2:D:1:MET:CA	2:D:2:ASN:HB3	2.51	0.40
2:D:164:TYR:O	2:D:165:LEU:HD22	2.22	0.40
2:D:1:MET:H2	2:D:2:ASN:HB3	1.84	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	175/190 (92%)	171 (98%)	4 (2%)	0	100	100
1	C	183/190 (96%)	173 (94%)	10 (6%)	0	100	100
2	B	209/218 (96%)	203 (97%)	6 (3%)	0	100	100
2	D	214/218 (98%)	208 (97%)	5 (2%)	1 (0%)	29	48
All	All	781/816 (96%)	755 (97%)	25 (3%)	1 (0%)	51	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	204	VAL

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	143/167 (86%)	136 (95%)	7 (5%)	25	47
1	C	147/167 (88%)	143 (97%)	4 (3%)	44	71
2	B	175/194 (90%)	173 (99%)	2 (1%)	73	89
2	D	176/194 (91%)	171 (97%)	5 (3%)	43	70
All	All	641/722 (89%)	623 (97%)	18 (3%)	43	70

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LYS
1	A	103	MET
1	A	123	LEU
1	A	133	ARG
1	A	156	PHE
1	A	160	GLN
1	A	190	ARG
2	B	16	GLN
2	B	208	MET
1	C	86	LEU
1	C	123	LEU
1	C	143	LYS
1	C	156	PHE
2	D	1	MET
2	D	88	ARG
2	D	134	GLU
2	D	138	GLN
2	D	163	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	ASN
2	B	104	GLN
2	D	31	HIS
2	D	138	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	181/190 (95%)	0.60	14 (7%) 13 13	43, 84, 134, 153	0
1	C	185/190 (97%)	0.48	14 (7%) 13 14	40, 68, 116, 145	0
2	B	211/218 (96%)	0.38	11 (5%) 27 29	42, 72, 125, 213	0
2	D	216/218 (99%)	0.52	23 (10%) 6 5	36, 75, 130, 156	0
All	All	793/816 (97%)	0.49	62 (7%) 13 13	36, 75, 130, 213	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	215	THR	6.5
1	C	145	GLY	6.2
2	D	212	ILE	6.0
2	D	216	LEU	5.8
1	A	88	THR	5.6
1	A	92	ASP	5.0
2	D	214	ASP	4.7
2	B	210	ARG	4.6
2	D	1	MET	4.6
2	B	1	MET	4.5
2	B	205	THR	4.3
2	B	2	ASN	4.3
1	A	160	GLN	4.2
1	C	91	GLY	3.8
1	A	162	SER	3.8
2	D	132	ALA	3.7
2	D	209	ARG	3.6
1	C	111	ILE	3.5
2	B	211	LEU	3.5
2	D	207	THR	3.5
1	C	159	VAL	3.5

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Mol	Chain	Res	Type	RSRZ
2	D	112	PHE	3.4
1	A	90	GLU	3.3
2	D	213	LYS	3.2
1	C	146	HIS	3.2
2	D	211	LEU	3.2
2	B	112	PHE	3.2
2	D	93	ASP	3.2
1	C	148	TYR	3.1
1	A	149	VAL	3.1
1	C	90	GLU	3.1
2	D	2	ASN	3.0
1	C	180	VAL	3.0
2	D	106	LYS	3.0
1	A	89	SER	3.0
1	A	91	GLY	2.9
2	B	115	GLU	2.9
1	A	7	SER	2.8
1	C	76	VAL	2.8
2	D	83	PHE	2.7
1	C	88	THR	2.6
1	C	107	CYS	2.6
2	B	204	VAL	2.6
1	A	111	ILE	2.5
1	C	89	SER	2.5
2	D	115	GLU	2.5
2	D	210	ARG	2.5
1	A	148	TYR	2.5
2	D	208	MET	2.5
2	B	30	LEU	2.4
1	A	165	GLY	2.4
2	B	94	GLY	2.3
2	B	83	PHE	2.3
2	D	133	THR	2.2
1	A	161	LEU	2.2
2	D	163	GLU	2.2
1	A	27	VAL	2.2
2	D	199	ALA	2.1
1	C	128	LEU	2.1
2	D	131	LEU	2.0
1	C	130	ALA	2.0
2	D	162	HIS	2.0



## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.